

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042820\
 Data File : BN010627.D
 Acq On : 28 Apr 2020 17:00
 Operator : CG/JU
 Sample : L2363-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 KY038PS038-200420

Quant Time: Apr 28 17:30:54 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN041320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 28 11:13:37 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	3035	0.40	ng	0.00
7) Naphthalene-d8	10.34	136	12904	0.40	ng	0.00
13) Acenaphthene-d10	14.21	164	8474	0.40	ng	0.00
19) Phenanthrene-d10	16.96	188	19899	0.40	ng	0.00
27) Chrysene-d12	21.16	240	17395	0.40	ng	0.00
34) Perylene-d12	23.33	264	15996	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.23	112	1095	0.16	ng	0.00
5) Phenol-d6	6.78	99	801	0.09	ng	0.00
8) Nitrobenzene-d5	8.72	82	3481	0.39	ng	0.00
11) 2-Methylnaphthalene-d10	11.94	152	7550	0.36	ng	0.00
14) 2,4,6-Tribromophenol	15.72	330	1646	0.41	ng	0.00
15) 2-Fluorobiphenyl	12.83	172	12457	0.40	ng	0.00
25) Fluoranthene-d10	19.00	212	22056	0.37	ng	0.00
29) Terphenyl-d14	19.61	244	19438	0.49	ng	0.00
Target Compounds						Qvalue
32) Bis(2-ethylhexyl)phthalate	21.10	149	1846	0.063	ng	95
33) Indeno(1,2,3-cd)pyrene	25.48	276	1024	0.020	ng	95
35) Benzo(b)fluoranthene	22.69	252	1070	0.020	ng	# 1
39) Benzo(g,h,i)perylene	26.12	276	913	0.021	ng	# 32

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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